### Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

# Listing of Claims:

- 1.-6. (Cancelled)
- 7. (Currently Amended) The A compound of Claim 4 wherein of formula (Ia):

#### wherein:

the A, C or D ring is independently fully saturated;

C1, C4, C11, C12, C15 and C16 are each independently substituted with two hydrogens;

C9 and C14 are each independently substituted with hydrogen:

C9 and C14 are each independently substituted with hydrogen,

R<sup>1</sup> is -OR<sup>7</sup>;

R2 is -R8-OR7;

R3 is -R10-N(R7)2:

R<sup>4a</sup> and R<sup>4b</sup> are each independently selected from hydrogen, alkenyl or alkynyl; or R<sup>4a</sup> is hydrogen,alkenyl or alkynyl and R<sup>4b</sup> is a direct bond to the carbon at C16; or R<sup>4a</sup> and R<sup>4b</sup> together form alkylidene or haloalkylidene;

R<sup>5</sup> is alkyl-or R<sup>5</sup> is a direct bond to the carbon at C14:

R<sup>6</sup> is hydrogen, -R<sup>8</sup>-OR<sup>7</sup> or -R<sup>8</sup>-N(R<sup>7</sup>)<sub>6</sub>:

each  $R^7$  is independently selected from the group consisting of hydrogen,  $-R^{10}$ -OR $^9$ ,  $-R^{10}$ -N( $R^9$ )<sub>2</sub>, alkyl, optionally substituted cycloalkyl, optionally substituted cycloalkyl, optionally substituted aryl, optionally substituted aralkyl, optionally substituted heterocyclylalkyl, optionally substituted heteroaryl and optionally substituted heteroarylalkyl;

each R<sup>8</sup> is independently selected from the group consisting of a direct bond, a straight or branched alkylene chain, and a straight or branched alkenylene chain;

each  $R^{\theta}$  is independently selected from the group consisting of hydrogen, alkyl, aryl and aralkyl; and

each  $R^{10}$  is independently selected from the group consisting of a straight or branched alkylene and a straight or branched alkylene chain,

<u>as a single stereoisomer, a mixture of stereoisomers, or as a racemic mixture of stereoisomers;</u>

or a pharmaceutically acceptable salt, solvate or prodrug thereof.

- (Currently Amended) The compound of Claim 7 selected from the group consisting of the following:
- $\label{eq:continuity} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-ethylideneoctahydroindene, ammonium chloride salt;$
- $\label{eq:continuity} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:controller} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium chloride salt;$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-aminoethyl)-7a\beta-methyl-1-methyleneoctahydroindene;$
- $\label{eq:continuity} 5-(1\beta\text{-methyl-}2\beta,4\beta\text{-dihydroxycyclohexyl})-4\alpha\text{-}(2\text{-aminoethyl})-7a\beta\text{-methyl-}1-\\ \text{methyleneoctahydroindene, ammonium acetate salt;}$
- $\label{eq:continuity} 5-(1\beta\text{-methyl-}2\beta,4\beta\text{-dihydroxycyclohexyl})-4\alpha\text{-aminomethyl-}7\alpha\beta\text{-methyl-}1-\\ methyleneoctahydroindene;}$
- $5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:control} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-difluoromethyleneoctahydroindene;$
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-aminomethyl-7aβ-methyl-1-difluoromethyleneoctahydroindene, ammonium chloride salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-aminomethyl-7aβ-methyl-1-dichloromethyleneoctahydroindene, ammonium chloride salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-aminomethyl-7aβ-methyl-1β-(propen-2-yl)ectahydroindene;

- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl) 4α-aminomethyl-7aβ-methyl-1β-(propen-2-yl)octahydroindene, ammonium acetate salt;
- $5-(1\beta-methyl-4\alpha,5\alpha-dihydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(4-dimethylaminobut-2Z-en-1-yl)-7a\beta-methyl-1-methyleneoctahydroindene;$
- 5-(1β-methyl-2β,4β-dihydroxycyclohexyl)-4α-(4-dimethylaminobut-2Z-en-1-yl)-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(ethyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt:
- 5-(1 $\beta$ -methyl-4 $\beta$ -hydroxy-2 $\beta$ -hydroxymethylcyclohexyl)-4 $\alpha$ -(benzyl)aminomethyl-7a $\beta$ -methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(cyclopropylmethyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt:
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(dimethyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene;
- $\label{eq:controller} 5-(1\beta-\text{methyl-}4\beta-\text{hydroxy-}2\beta-\text{hydroxymethylcyclohexyl)-}4\alpha-(\text{dimethyl})\text{-}\text{aminomethyl-}7\alpha\beta-\text{methyl-}1-\text{methyl-}\text{neoctahydroindene, ammonium acetate salt;}$
- $\label{eq:controller} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-(methyl) aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene;$
- 5-(1 $\beta$ -methyl-4 $\beta$ -hydroxy-2 $\beta$ -hydroxymethylcyclohexyl)-4 $\alpha$ -(methyl)aminomethyl-7a $\beta$ -methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- $\label{eq:controller} 5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-(2-methylpropyl) aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene;$
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(2-methylpropyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(1-methylpiperidin-4yl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium diacetate salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(3-nitrobenzyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(piperonyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt;

- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(pyrrol-2ylmethyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt:
- $5-(1\beta-methyl-4\beta-hydroxy-2\beta-hydroxymethylcyclohexyl)-4\alpha-(furfuryl)aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate sait;$
- 5-(1β-methyl-4β-hydroxy-2β-hydroxymethylcyclohexyl)-4α-(pyridin-3ylmethyl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate sait;
- $5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-methylpropyl)$ aminomethyl- $7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt:$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(pyridin-3-ylmethyl) aminomethyl-7a\beta-methyl-1-methylene octahydroindene:$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-hydroxyethyl) aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- 5-(1β-methyl-2β,4β-dihydroxycyclohexyl)-4α-(furfuryl)aminomethyl-7aβ-methyl-1-methyleneoctahydroindene, ammonium acetate salt;
- $\label{eq:controller} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-dimethylaminoethyl)aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:condition} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-cyclohex-1-en-1-ylethyl)aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(2-morpholin-4-ylethyl) aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene, ammonium acetate salt;$
- $\label{eq:continuity} 5-(1\beta-methyl-2\beta,4\beta-dihydroxycyclohexyl)-4\alpha-(3-methylphenyl) aminomethyl-7a\beta-methyl-1-methyleneoctahydroindene;$
- $\label{eq:continuity} 5-(1\beta-\text{methyl-}2\beta,4\beta-\text{dihydroxycyclohexyl})-4\alpha-(\text{benzyl})\\ a minomethyl-7a\beta-\text{methyl-1-methyleneoctahydroindene}; and$
- $\label{eq:continuity} 5-(1\beta-\text{methyl-}2\beta,4\beta-\text{dihydroxycyclohexyl})-4\alpha-(2-(3-\text{methyl-henyl})a\text{minoethyl})-7a\beta-\text{methyl-1-methyleneoctahydroindene}.$

#### 9.-21 (Cancelled)

22. (Currently Amended) A pharmaceutical composition comprising a pharmaceutically acceptable excipient and a compound of formula (la):

### wherein:

the A, C or D ring is independently fully saturated;

C1, C4, C11, C12, C15 and C16 are each independently substituted with two hydrogens;

C9 and C14 are each independently substituted with hydrogen:

 $R^1$  is  $-OR^7$ ;

R2 is -R8-OR7;

 $R^3$  is  $-R^{10}-N(R^7)_2$ ;

R<sup>4a</sup> and R<sup>4b</sup> together form alkylidene or haloalkylidene;

R<sup>5</sup> is alkyl;

R6 is hydrogen, -R8-OR7 or -R8-N(R7)2;

each R<sup>7</sup> is independently selected from the group consisting of hydrogen, -R<sup>10</sup>-OR<sup>9</sup>, -R<sup>10</sup>-N(R<sup>9</sup>)<sub>2</sub>, alkyl, optionally substituted cycloalkyl, optionally substituted cycloalkyl, optionally substituted arriv, optionally substituted arriv, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl;

each  $R^8$  is independently selected from the group consisting of a direct bond, a straight or branched alkylene chain, and a straight or branched alkylene chain;

 $\underline{\text{each } R^9 \text{ is independently selected from the group consisting of hydrogen, alkyl, anyl and}} \\ \text{aralkyl: and}$ 

 $\underline{\text{each } R^{10} \text{ is independently selected from the group consisting of a straight or branched}} \\ \underline{\text{alkylene and a straight or branched alkenylene chain.}}$ 

as a single stereoisomer, a mixture of stereoisomers, or as a racemic mixture of stereoisomers;

or a pharmaceutically acceptable salt, solvate or prodrug thereofcompound of Claim (I):

wherein:

the A, C or D ring is independently fully saturated, partially saturated or fully unsaturated; C1, C4, C11, C12, C15 and C16 are each independently substituted with two of the following, which are independently selected: hydrogen, alkyl, -R\*-OR\*, or -R\*-N(R\*)<sub>2r</sub> provided that C4 is not substituted by two methyl groups;

C9 and C14 are each independently-substituted with hydrogen, alkyl,  $-R^8$ -OR $^2$ , or  $-R^8$ -N/R $^2$  $)_{22}$ :

R1-is-OR7-or-N(R7)2:

R<sup>2</sup>-and R<sup>3</sup>-are each independently selected from the group-consisting of —R<sup>8</sup>-OR<sup>7</sup>,

—R<sup>8</sup>-OC(O)R<sup>9</sup>,—R<sup>10</sup>-N(R<sup>2</sup>)<sub>a</sub>.—R<sup>10</sup>-N(R<sup>9</sup>)C(O)R<sup>9</sup>,—R<sup>10</sup>-N(R<sup>9</sup>)S(O)<sub>A</sub>R<sup>2</sup> (where t is 1 or 2),

—R<sup>10</sup>-N(R<sup>9</sup>)C(NR<sup>9</sup>)N(R<sup>9</sup>)<sub>a</sub>, alkyl, alkenyl, optionally substituted aralkyl, optionally substituted aralkyl, optionally substituted heteroarylalkyl, optionally substituted heteroarylalkyl, optionally substituted heteroarylalkonyl, and optionally substituted heteroarylalkonyl;

R<sup>4a</sup>-and R<sup>4b</sup>-are-each independently-selected from hydrogen, alkenyl-or-alkynyl; or R<sup>4a</sup>-is-hydrogen, alkenyl-or-alkynyl-and R<sup>4b</sup>-is-a direct bond to the carbon at C16; or R<sup>4a</sup>-and R<sup>4b</sup>-together form alkylidene or haloalkylidene;

R<sup>5</sup>-is alkyl or R<sup>5</sup>-is a direct bond to the carbon at C14;

R<sup>6</sup> is hydrogen, -R<sup>8</sup>-OR<sup>2</sup> or -R<sup>8</sup>-N(R<sup>2</sup>)2;

each R<sup>7</sup>-is independently selected from the group consisting of hydrogen,—R<sup>10</sup>-OR<sup>9</sup>, -R<sup>10</sup>-N(R<sup>9</sup>)<sub>2n</sub>-alkyl, optionally substituted cycloalkyl, optionally substituted cycloalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl;

each R<sup>8</sup> is independently selected from the group-consisting of a direct bond, a straight or branched alkylene chain, and a straight or branched alkenylene chain;

each  $\mathbb{R}^{\theta}$  is independently selected from the group consisting of hydrogen, alkyl, aryl and aralkyl; and

each R<sup>40</sup> is independently selected from the group consisting of a straight or branched alkylene and a straight or branched alkylene chain;

as a single stereoisomer, a mixture of stereoisomers, or as a racemic mixture of stereoisomers:

or a pharmaceutically acceptable salt, solvate or prodrug thereof.

# 23. (Cancelled)

24. (Withdrawn and Currently Amended) A method of treating an inflammatory condition or disease in a mammal, which method comprises administering to the mammal in need thereof a therapeutically effective amount of a compound having the following formula (Ia):

#### wherein:

the A, C or D ring is independently fully saturated;

C1, C4, C11, C12, C15 and C16 are each independently substituted with two hydrogens;

C9 and C14 are each independently substituted with hydrogen;

 $R^1$  is  $-OR^7$ ;

R2 is -R8-OR7;

 $R^3$  is  $-R^{10}-N(R^7)_2$ ;

R<sup>4a</sup> and R<sup>4b</sup> together form alkylidene or haloalkylidene;

R<sup>5</sup> is alkyl;

R<sup>6</sup> is hydrogen, -R<sup>8</sup>-OR<sup>7</sup> or -R<sup>8</sup>-N(R<sup>7</sup>)<sub>2</sub>;

each R<sup>7</sup> is independently selected from the group consisting of hydrogen, -R<sup>10</sup>-OR<sup>9</sup>,
-R<sup>10</sup>-N(R<sup>9</sup>)<sub>2</sub>, alkyl, optionally substituted cycloalkyl, optionally substituted cycloalkyl,
optionally substituted aryl, optionally substituted arelkyl, optionally substituted heterocyclylalkyl,

optionally substituted heteroaryl and optionally substituted heteroarylalkyl;

each R<sup>8</sup> is independently selected from the group consisting of a direct bond, a straight or branched alkylene chain, and a straight or branched alkylene chain:

 $\underline{\text{each } R^9 \text{ is independently selected from the group consisting of hydrogen, alkyl, anyl and} \\ \underline{\text{aralkyl}}; \text{and}$ 

each  $R^{10}$  is independently selected from the group consisting of a straight or branched alkylene and a straight or branched alkylene chain,

as a single stereoisomer, a mixture of stereoisomers, or as a racemic mixture of stereoisomers:

or a pharmaceutically acceptable salt, solvate or prodrug thereofcompound of formula (H):

wherein:

the A, C or D ring is independently fully saturated, partially saturated or fully unsaturated; C1, C4, C11, C12, C15 and C16 are each independently substituted with two of the following, which are independently selected: hydrogen, alkyl,—R\*-OR\*, or—R\*-N(R\*)<sub>2r</sub>, provided that C4 is not substituted by two methyl-groups;

C9 and C14 are each independently-substituted with hydrogen, alkyl, \_R\*-QR\*, or \_R\*-N(R\*);;

R<sup>2</sup> and R<sup>3</sup> are each independently-selected from the group consisting of —R<sup>8</sup>–OR<sup>2</sup>, -R<sup>8</sup>–OC(O)R<sup>9</sup>,—R<sup>10</sup>-N(R<sup>2</sup>)<sub>2</sub>,—R<sup>10</sup>-N(R<sup>9</sup>)C(O)R<sup>9</sup>,—R<sup>10</sup>-N(R<sup>9</sup>)S(O)<sub>2</sub>R<sup>2</sup> (where t is 1 or 2), -R<sup>10</sup>-N(R<sup>9</sup>)C(NR<sup>9</sup>)N(R<sup>9</sup>)<sub>2</sub>, alkyl, alkenyl, optionally substituted aralkyl, optionally substituted aralkyl, optionally substituted heteroarylalkyl, optionally substituted heteroarylalkyl, optionally substituted heteroarylalkenyl;

 $R^{4a}$  and  $R^{4b}$  are each independently selected from hydrogen, alkenyl-or-alkynyl; or  $R^{4a}$  is hydrogen, alkenyl-or-alkynyl and  $R^{4b}$  is a direct bond to the carbon at C16; or  $R^{4a}$  and  $R^{4b}$  -together-form alkylidene or haloalkylidene;

R<sup>5</sup> is alkyl or R<sup>5</sup> is a direct bond to the carbon at C14;

R<sup>6</sup> is hydrogen, R<sup>8</sup>-OR<sup>7</sup> or -R<sup>8</sup>-N(R<sup>7</sup>)2;

each R<sup>z</sup>-is-independently-selected from the group consisting of hydrogen, –R<sup>10</sup>-OR<sup>9</sup>, –R<sup>10</sup>-N(R<sup>9</sup>)<sub>zn</sub>-alkyl, optionally substituted cycloalkyl, optionally substituted cycloalkylalkyl, optionally substituted aryl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl, optionally substituted heterocyclylalkyl,

each R<sup>8</sup>-is independently selected from the group consisting of a direct bond, a straight or branched alkylene chain, and a straight or branched alkylene chain;

each  $\mathbb{R}^9$  is independently selected from the group consisting of hydrogen, alkyl, aryl and aralkyl; and

each R<sup>19</sup>-is independently selected from the group consisting of a straight or branched alkylene and a straight or branched alkylene shain:

as a single stereoisemer, a mixture of stereoisemers, or as a racemic mixture of stereoisemers:

or a pharmaceutically acceptable salt, solvate or prodrug thereof.

25. (Cancelled)

26. (Withdrawn and Currently Amended) The method of Claim 24 or Claim 25 wherein the inflammatory condition or disease is selected from the group consisting of the following:

arthritis (including rheumatoid arthritis, psoriatic arthritis, ankylosing spondylitis, osteoarthritis, gout, and synovitis), inflammations of the brain (including multiple sclerosis, Alzheimer's, AIDS dementia, stroke, encephalitis, trauma, and Creutzfeld-Jakob disease), inflammatory bowel disease (including Crohn's disease and ulcerative colitis), irritable bowel syndrome, ischemia-reperfusion injury (including myocardial infarction), sarcoidosis, psoriasis, tissue/organ transplant, graft vs host disease, systemic lupus erythematosus, Type I juvenile diabetes, vasculitis, artherosclerosis, cardiomyopathy, autoimmune myocarditis, atopic dermatitis, asthma, allergy, allergic rhinitis, and chronic obstructive pulmonary disease (including emphysema and bronchitis).

27.-29. (Cancelled)